

Current-spin coupling for ferromagnetic domain walls in fine wires

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The coupling between a current and a domain wall is examined. In the presence of a finite current and the absence of a potential which breaks the translational symmetry, there is a perfect transfer of angular momentum from the conduction electrons to the wall. As a result, the ground state is in uniform motion. This remains the case when relaxation is accounted for. This is described by, appropriately modified, Landau-Lifshitz-Gilbert equations.

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Spintronic devices have great technological promise but represent a challenging problem at both an applied and fundamental level. It has been shown theoretically[1, 2] that the direction of a magnetic domain might be switched using currents alone. Devices designed to use this principle often consist of multi-layers of magnetic and non-magnetic conductors. The advantages of similar devices based upon the current induced displacement of a *domain wall* are simplicity and the fact that the switching current is much smaller[3, 4, 5, 6]. Experimentally the current induced displacement of a domain wall has been clearly demonstrated and in recent experiments[5, 6] the velocity of the wall was measured.

The current induced motion of a magnetic domain involves the transfer of angular momentum from the conduction electrons. The early theory[1, 2] and most of the subsequent work[7] is based upon some type of assumption about this torque transfer process and there is no real consensus on how this should appear in the (Landau-Lifshitz-Gilbert) equations of motion[7]. The purpose of this Letter is to develop a *complete* first principles theory of this process for a *domain wall*, this based upon a specific model Hamiltonian and physically justified approximations. This is intended to form the basis for the more complicated problem of multi-layers.

Although similar conclusions are valid for the Stoner, and related, models, here attention will be focused upon the *s-d*-exchange model. The direction of the local moments, \vec{S}_i , which make up the domain wall are specified by the usual Euler angles θ_i and ϕ_i . To make diagonal the interaction $+J\vec{S}_i \cdot \vec{s}_i$, at site *i*, the axis of quantization of the conduction electrons, \vec{s}_i , is rotated along this same direction. If $\psi(\vec{r}_i, t)$ is the conduction electron spinor field then this amounts to making a SU(2) gauge transformation, $\psi(\vec{r}_i, t) \rightarrow r(\theta_i, \phi_i)\psi(\vec{r}_i, t)$, where $r(\theta_i, \phi_i) \equiv e^{i\phi_i s_z} e^{i\theta_i s_y} e^{i\phi_i s_z} = (\cos(\theta_i/2) + i \sin(\theta_i/2) \sin \phi_i \sigma_x - i \sin(\theta_i/2) \cos \phi_i \sigma_y)$ and where $\vec{\sigma}$ are the Pauli matrices. If there was SU(2) gauge *invariance*, there would exist *three* gauge particles, the analogy of the W^\pm and Z which mediate the weak interaction. However here, without this invariance, this transformation still introduces *three* gauge fields. The *longitudinal* such field has been exploited in the development of theories of the Hall effect[8].

It is Bazaliy et al.[9], who first developed a theory of the torque transfer to a magnetic domain, for the half-metal limit, based upon this same field. Their formalism generates a *transverse* field in the Landau-Lifshitz equations through a finite derivative $\partial \mathcal{V}\{\theta_i, \phi_i\}/\partial \phi_i$ where $\mathcal{V}\{\theta_i, \phi_i\}$ is the expectation value of the energy as a function of the angles $\{\theta_i, \phi_i\}$. *However*, a finite ground state expectation value for such a derivative implies that the solution is not stable and it follows that their finite velocity solution must relax to one which is *stationary* with a finite common value of the $\phi_i = \phi_0$, this anticipating the conclusions of Tatara and Kohno[10].

Here it will be shown that the torque transfer effects can be correctly accounted for in a simpler U(1) theory in which the only generators are S_{iy} for rotations about the axis perpendicular to the wall plane. The local spins are usually co-planer, see below, and the rotations $r(\theta_i) \equiv e^{i\theta_i s_y} = (\cos(\theta_i/2) - i \sin(\theta_i/2) \sigma_y)$ are all that is needed to diagonalize $+J\vec{S} \cdot \vec{s}$. This simpler approach can *only* generate a *transverse* gauge field, i.e., the one ignored in the earlier work[9], and which *alone* is found to be the origin of the torque transfer process. The longitudinal field, along with $\partial \mathcal{V}\{\theta_i, \phi_i\}/\partial \phi_i$, and the ground state ϕ_0 are strictly null, and the correctly *relaxed* ground state *does* have a finite velocity in the absence of pinning.

The *s-d*-exchange Hamiltonian is,

$$\begin{aligned} \mathcal{H} = & - \sum_{\langle ij \rangle \sigma \sigma'} \left(c_{i\sigma}^\dagger t_{ij\sigma\sigma'} c_{j\sigma'} + H.c. \right) - \mu \hat{N} \\ & - \sum_i (J \vec{S}_i \cdot \vec{s}_i + A^0 S_{iz}^2 - K_\perp^0 S_{iy}^2) \\ & - J_s^0 \sum_{\langle ij \rangle} \vec{S}_i \cdot \vec{S}_j \end{aligned} \quad (1)$$

where $c_{i\sigma}^\dagger$ is the conduction electron creation operator for spin σ and site *i*. The uniform hopping integral is $t_{ij\sigma\sigma'} = t \delta_{\sigma\sigma'}$ (the strange notation is for later use), and \hat{N} is number operator. The exchange constants are *J* and J_s^0 for the coupling between the local moments and the conduction electrons and for the direct coupling between local moments. With the anisotropy constant $A^0 > 0$ the *z* direction is an easy axis while, with $K_\perp^0 > 0$, the *y* axis is hard. These anisotropy constants include both

intrinsic anisotropy plus the effects of a demagnetizing field.

It is first necessary to describe the domain wall. The Holstein-Primakoff transformation[11] $S_{iz} = S - b_i^\dagger b_i$, $S_i^+ = (2S - b_i^\dagger b_i)^{1/2} b_i \approx (2S)^{1/2} b_i$ is used to quantize the spin degrees of freedom using the direction defined by θ_i and ϕ_i as the axis of quantization for the local spins. To within a constant, for the spin Hamiltonian in Eqn. (1), $\mathcal{V}\{\theta_i, \phi_i\} = -AS^2 \cos^2 \theta_i + K_\perp S^2 \sin^2 \theta_i \sin^2 \phi_i - \sum_{\langle ij \rangle} J_s \cos \theta_{ij}$ where $A = [(2S - 1)/2S]A^0$, $K_\perp = [(2S - 1)/2S]K_\perp^0$, and θ_{ij} is the angle between the spins at sites i and j . The energy $K_\perp S^2 \sin^2 \theta_i \sin^2 \phi_i$ favors a wall with $\phi_i = 0$, whence assuming a slow variation in the angle θ_i , i.e., that $\theta_{ij} \approx \nabla \theta_i \cdot \vec{r}_{ij}$, where $\vec{r}_{ij} = \vec{r}_j - \vec{r}_i$ is the vector which joins near neighbor planes of the wall in the z direction ($r_{ij} = a$), and again dropping a constant $\mathcal{V}\{\theta_i\} = +J_s S^2 (\nabla \theta_i \cdot \vec{r}_{ij})^2 - AS^2 \cos^2 \theta_i$. The domain wall structure, which lies in the $x - z$ -plane, can be determined by minimizing this $\mathcal{V}\{\theta_i\}$, and the standard result is $\theta(z_i) = 2 \cot^{-1} e^{-(z_i/w)}$, where z_i is coordinate of site i . The wall width $w = a(J_s/2A)^{1/2}$. The local spins rotate so $\theta = 0$ becomes $\theta = \pi$ with increasing z .

In order to have precise expressions for the conjugate position and momentum of the wall, it is observed that the later is the generator of displacements. The product of small y -axis rotations $R(\Delta z) = \prod_i \exp(i(\theta_i - \theta_j)(\Delta z/a)(S_{iy}/\hbar)) \approx (1 + (1/2) \sum_i (\nabla \theta_i \cdot \vec{r}_{ij}) \Delta(z/a)(2S)^{1/2}(b_i^\dagger - b_i))$, produces a wall translation by Δz , and so the momentum, per atom in a plane of the wall, $p_z = i(\hbar/2)(v_c/a^2 \mathcal{A}) \sum_i (\nabla \theta_i \cdot \vec{r}_{ij})(2S)^{1/2}(b_i^\dagger - b_i)$, where v_c is the cell volume and \mathcal{A} is the cross-sectional area of the wire. The Goldstone boson which restores the translational invariance of \mathcal{H} can then be identified as,

$$b_w^\dagger = (J/8A)^{1/4}(v_c/a\mathcal{A}) \sum_i (\nabla \theta_i \cdot \vec{r}_{ij}) b_i^\dagger,$$

where the constant of proportionality reflects the requirement that $[b_w, b_w^\dagger] = 1$ and uses the explicit wall solution for $\theta(z)$. (The modes which disperse from this Goldstone boson are excitations of the wall.) Thus $\hat{p}_z = i\hbar(2A/J)^{1/4}(2S)^{1/2}(1/a)(b_w^\dagger - b_w)$ while the conjugate coordinate, such that $[\hat{z}_0, \hat{p}_z] = i\hbar$, is $\hat{z}_0 = a(J/2A)^{1/4}(S/2)^{-1/2}(b_w^\dagger + b_w)$.

Returning to the description of the conduction electrons, performing the above described SU(2) rotations, the conduction electron part of Eqn. (1) reduces to,

$$\mathcal{H}_e = - \sum_{\langle ij \rangle > \sigma \sigma'} \left(c_{i\sigma}^\dagger t_{ij\sigma\sigma'} c_{j\sigma'} + H.c. \right) - JS \sum_i s_{iz} - \mu \hat{N}, \quad (2)$$

in which now $t_{ij} = tr^{-1}(\theta_i, \phi_i)r(\theta_j, \phi_j)$. But since $\phi_i = 0$ this simplifies to the U(1) result:

$$t_{ij} = t(\cos(\theta_i/2) + i \sin(\theta_i/2)\sigma_y)(\cos(\theta_j/2) - i \sin(\theta_j/2)\sigma_y). \quad (3)$$

which is, $t_{ij} = t[\cos(\theta_i/2)\cos(\theta_j/2) + \sin(\theta_i/2)\sin(\theta_j/2) + i(\sin(\theta_i/2)\cos(\theta_j/2) - \sin(\theta_j/2)\cos(\theta_i/2)\sigma_y]$, using $\sigma_y^2 = 1$. The standard identities then lead to $t_{ij} = t[\cos \nabla \theta_i \cdot \vec{r}_{ij} + i(\nabla \theta_i \cdot \vec{r}_{ij})\sigma_y]$, and correct to second order in the gradient, this can be written as,

$$t_{ij} = e^{i \int_{\vec{r}_i}^{\vec{r}_j} \vec{A} \cdot d\vec{r}} t \cos(\nabla \theta_i \cdot \vec{r}_{ij}); \quad \vec{A} = (\nabla \theta_i \cdot \vec{r}_{ij}) s_y \hat{z} \quad (4)$$

In order that this central result be valid, the conduction electrons must follow the local spin magnetization as they pass through the wall and this requires that the adiabatic theorem be satisfied. When the transverse field represented by $\Delta t_{ij}^\perp = it(\nabla \theta_i \cdot \vec{r}_{ij})\sigma_y$ is ignored (the equivalent to $\vec{A} = 0$), the eigenstates of \mathcal{H}_e are also eigenstates of the total angular momentum $\sum_i (\vec{S}_i + \vec{s}_i)$ evaluated in the *local* frame. Given that $J > 0$, in the ground state, \vec{S}_i and \vec{s}_i are then parallel.

This approximation is manifestly valid in the half metal limit when $J \gg t$. The ground state is then a mixture of states in which all sites are either singly occupied by an electron or unoccupied. The singly occupied sites with the maximum angular momentum $S + (1/2)$ have the lowest energy while other states, and those with two electrons per site, have an energy which is higher by $\sim J$ and hence have negligible weight in the ground state. The Wigner-Eckart theorem then dictates that all the matrix elements of \vec{s}_i are equal to those of $(\vec{S}_i/2S)$.

However, since spatially the magnetization rotates slowly, a much weaker inequality suffices. The adiabatic theorem simply demands that the transverse field $\Delta t_{ij}^\perp = it(\nabla \theta_i \cdot \vec{r}_{ij})\sigma_y$ be small compared to the longitudinal field Jm_s . The wall rotates by π over a distance w so $\nabla \theta_i \sim \pi/w$ and $\Delta t_{ij}^\perp \sim i\pi t(a/w)$, and required is,

$$\pi t(a/w) \ll Jm_s \quad (5)$$

which since, e.g., for Permalloy $w/a \sim 10^3$ is typically well satisfied. The conduction electron magnetization comprises two components with, by definition, the (minority) majority conduction electrons (anti-) parallel to the axis of quantization, i.e., the direction of the local spin. In the *local frame* the majority (minority) electrons have $\sigma_z = +1$ ($\sigma_z = -1$), so that it follows that when the adiabatic theorem is satisfied,

$$\vec{s}_i = \sigma_z (\vec{S}_i/2S), \quad (6)$$

independent of the details of the electronic structure, etc.

The non-interacting approximation is therefore to indeed ignore Δt_{ij}^\perp . If it was not for the cosine in $t_{ij} = t \cos(\nabla \theta_i \cdot \vec{r}_{ij})$ the problem is then identical to that without the wall. Using again $(\nabla \theta_i \cdot \vec{r}_{ij}) \sim \pi a/w$, the correction is $\sim t\pi^2(a/w)^2$ which with $(a/w) \sim 10^{-3}$ might be safely ignored, i.e., at this level the electronic structure is unchanged by the wall and since the reflection probability is small, there is a negligible *pressure* exerted by the conduction electrons. However, in the spin sector,

this correction *is* important. Evaluating the coefficient of $(\nabla\theta_i \cdot \vec{r}_{ij})^2$ in first order perturbation theory leads to, the renormalization $J_s = J_s^0 + (x't/2S^2)$, of the exchange coupling[13]. The effective concentration $x' = \langle c_i^\dagger c_j \rangle$.

At this lowest level of approximation the wall is stationary and is put in motion only when the torque transfer Δt_{ij}^\perp -term is accounted for as a perturbation. Even with this term, the effective Hamiltonian, Eqn. (2), is evidently of single particle nature. Consider first the single particle description of the majority spin electrons. In order to account for the torque transfer term proportional to Δt_{ij}^\perp , use is made of Eqn. (6). To this end, it is noted, for majority electrons $\sum_{\sigma\sigma'} c_{i\sigma}^\dagger \sigma_{y\sigma\sigma'} c_{j\sigma'} = i(s_i^- c_{i\uparrow}^\dagger c_{j\uparrow} - c_{i\uparrow}^\dagger c_{j\uparrow} s_j^+)$. Then by virtue of Eqn. (6), e.g., $s_i^+ \approx (2S)^{-1/2} b_i$, and $\sum_{\sigma\sigma'} c_{i\sigma}^\dagger \sigma_{y\sigma\sigma'} c_{j\sigma'} = i(2S)^{-1/2} c_{i\uparrow}^\dagger c_{j\uparrow} (b_i^\dagger - b_j)$. Combining this with the similar result for the minority electrons, the Δt_{ij}^\perp -term becomes,

$$\mathcal{H}_\tau = -\frac{t}{2(2S)^{1/2}} \sum_{ij\sigma} \sigma (\nabla\theta_i \cdot \vec{r}_{ij}) c_{i\sigma}^\dagger c_{j\sigma} (b_i - b_j^\dagger) + H.c., \quad (7)$$

which couples the spin current to the magnons and reflects the *entire* torque transfer process. The necessary correction is obtained, directly, by taking the expectation value with respect to the conduction electrons, whence Eqn. (7) reduces to

$$\mathcal{H}_\tau = -i \frac{\hbar j_s a^2}{2eS} \sum_i (\nabla\theta_i \cdot \vec{r}_{ij}) (2S)^{1/2} (b_i - b_i^\dagger), \quad (8)$$

where j_s is the *spin* current. Here b_{i+1}^\dagger is replaced by b_i^\dagger , an approximation valid in the continuum limit $w \gg a$. The quantity $(2S)^{1/2} (b_i - b_i^\dagger) \approx S_{iy}$ and \mathcal{H}_τ corresponds to an effective field which is *strictly* transverse. The appearance of such a term linear in $(b_i - b_i^\dagger)$ signals that there is no time independent solution.

Comparing \mathcal{H}_τ with the definition of \hat{p}_z makes evident that $\mathcal{H}_\tau \propto j_s \hat{p}_z$, i.e., the current *only* couples to the collective coordinate of the wall via its momentum. The effect of \mathcal{H}_τ is then to put the wall in motion along the z -direction, and it is necessary to study the time dependent Schrödinger equation: $i\hbar(\partial/\partial t)\psi(\vec{r}_i, t) = \mathcal{H}\psi(\vec{r}_i, t)$. The effect of adding the rotations, r , which displace the wall is $i\hbar(\partial/\partial t) \rightarrow i\hbar r(\partial/\partial t)r^{-1} = i\hbar(\partial/\partial t) - \hbar(\partial\theta_i/\partial t)M_{iy}$, where $\vec{M}_i = \vec{S}_i + \vec{s}_i$ is the *total* spin angular momentum. This generates a second purely transverse field term in the effective Hamiltonian:

$$\hbar \sum_i \frac{\partial\theta_i}{\partial t} M_{iy} = \hbar \frac{M}{S} \sum_i \frac{\partial\theta_i}{\partial t} (2S)^{1/2} (b_i - b_i^\dagger)$$

using the fact that $(\vec{M}_i/M) = (\vec{S}_i/S)$, i.e., that all magnetizations are parallel. Thus when

$$\frac{\partial\theta_i}{\partial t} = v_0 \frac{\partial\theta_i}{\partial z}; \quad v_0 = \frac{j_s a^3}{2M}. \quad (9)$$

the effective fields generated by the spatial and temporal rotations of the axes of quantization cancel each other. Given $\theta_i = \theta(z_i) \equiv 2 \cot^{-1} e^{-(z_i/w)}$ for $j_s = 0$, the new *ground state* has $\theta_i = \theta(z_i - v_0 t)$, i.e., the wall moves without distortion and without tilting or twisting. It is easy to show that the result $v_0 = (j_s v_c / 2M)$ reflects the conservation of the z -component of the total angular momentum, i.e., that the net spin current, $2j_s$ carried towards the wall by the electrons equals the change in the angular momentum, $j_s = Mv/v_c$, of the wall due to its motion. Since the conduction electrons are polarized $j_s = pj$ is related to the charge current j by some material determined parameter p .

It is to be observed that by making these specific time dependent rotations the *net* transverse magnetic field has vanished from the problem. Thus the conditions for equilibrium in rotating frame are *identical* to those for the static problem, when $j_s = 0$, and in particular $\partial\mathcal{V}/\partial\phi_i = 0$ since this is an equilibrium condition. In order to verify this point of divergence with Bazaliy et al.[9], it is useful to calculate this derivative directly. This requires the full SU(2) transformations. After some algebra the result is, Eqn. (4) with, $\vec{A} = (1/2)(\nabla\theta_i)\sigma_y' + (1/2)\sin\theta_i(\nabla\phi_i)\sigma_x' + (1 - \cos\theta_i)(\nabla\phi_i)\sigma_z'$ where the σ' are defined in the local frame of reference. Directly, both new terms are zero for a simple domain wall for which $\nabla\phi_i = 0$, and $\partial\mathcal{V}/\partial\phi_i = 0$ [14].

In terms of the Landau-Lifshitz equations, in the laboratory frame,

$$\frac{D\vec{M}}{Dt} \equiv \frac{\partial\vec{M}}{\partial t} - (\vec{j}_s \cdot \nabla)\vec{M} = g\mu_B \vec{M} \times \vec{B} \quad (10)$$

where the effective field is ($K_\perp = 0$ for simplicity):

$$\vec{B} \equiv \frac{\partial\mathcal{V}}{\partial\vec{M}} = JS^2 a^2 \nabla^2 \vec{M} + \frac{2A}{M^2} (\hat{z} \cdot \vec{M}) \hat{z}. \quad (11)$$

This is formally similar to the result obtained by Bazaliy et al.[9], in the half metal limit, *except* for the fact that the term proportional to j_s does *not* arise from a finite $\partial\mathcal{V}/\partial\phi$, via Eqn. (11). Also this torque transfer term arises from Δt_{ij}^\perp which is simply *not* included in that work. Equation (10) defines the “particle derivative” $D\vec{M}/Dt$, i.e., that at a fixed position in the *moving* wall.

This different origin of the torque transfer term is of particular importance when relaxation is accounted for. With $j_s = 0$, relaxation is traditionally included through a Gilbert term $-(\alpha/M)\vec{M} \times (\partial\vec{M}/\partial t)$, parameterized by α , and which assures the system relaxes until \vec{M} is parallel to the internal field \vec{B} , and which corresponds to an *absolute minimum* of $\mathcal{V}\{\theta_i, \phi_i\}$. This is an evident requirement of the *second law of thermodynamics*. In order that this law be satisfied when j_s is finite, the Gilbert term must involve rather $D\vec{M}/Dt$ and the Landau-Lifshitz-Gilbert equations are,

$$\frac{D\vec{M}}{Dt} = g\mu_B \vec{M} \times \vec{B} - \frac{\alpha}{M} \vec{M} \times \frac{D\vec{M}}{Dt}. \quad (12)$$

with the present Eqn. (11).

The justification, at the microscopic level, of this form of relaxation is complicated. For the micron sized system it is widely assumed that the principal loss mechanism corresponds spin-orbit scattering of the conduction electrons[7]. Above a few tens of degrees, the Yafet-Elliot [15], spin-orbit scattering off phonons is almost certainly dominant. However, here there is little interest in obtaining a detailed expression for $(1/\tau)$ and so rather considered is the simplest model,

$$\mathcal{H}_{\text{so}} = i\lambda \sum_i \sum_{\vec{k}\vec{k}'} \sum_{\sigma\sigma'} e^{i(\vec{k}-\vec{k}')\cdot\vec{r}_i} \vec{\sigma}_{\sigma\sigma'} \cdot (\vec{k} \times \vec{k}') c_{\vec{k}\sigma}^\dagger c_{\vec{k}'\sigma'}$$

which represents such scattering off impurities, located are \vec{r}_i , and which enters the theory in the same manner as does phonon scattering. In fact, if the concentration c of “impurities” is taken to be $\sim T/T_D$ where T_D is the Deby temperature, this mimics the effect of phonons at modest temperatures. The problem is treated in the *rotating frame* since this reduces to zero the gauge fields created by the current j_s . Now in \mathcal{H}_{so} it is necessary to make the replacement $\vec{\sigma} \rightarrow r\vec{\sigma}r^{-1}$. The angular speed of rotation, $\sim \pi v_0/w \sim 10^3$ rad/s, is by far the lowest frequency in the problem and hence it can be assumed that, $\omega_0 = \dot{\theta}_i$ is a constant during the relaxation process. Further, in practice, the wall width w is large compared to the mean free path, and so the spatial dependence of ω_0 can be ignored. The problem is then very similar to the linear response to an applied radio frequency field. In the context of dilute magnetic alloys, this has been studied in some detail[16]. The Gilbert term is found to be of the correct form with, in the present notation, $\alpha = 1/(\omega_s\tau)$ where $\omega_s = g\mu_B B$ and, $(1/\tau) = W_{\uparrow\rightarrow\downarrow} - W_{\downarrow\rightarrow\uparrow}$, and where, e.g., $W_{\uparrow\rightarrow\downarrow}$ is the transition rate for conduction electrons to flip from up to down spin given essentially by the “Golden Rule”. Assuming a random concentration of “impurities”, e.g., $W_{\uparrow\rightarrow\downarrow} = c(\pi\lambda^2/\hbar) \sum_{\vec{k}\vec{k}'} [(k_z k'_x - k_x k'_z)^2 \delta(\epsilon_{\vec{k}} - \epsilon_{\vec{k}'}) + (1/4)((k_y k'_z - k_z k'_y)^2 + (k_x k'_y - k_y k'_x)^2) \delta(\epsilon_{\vec{k}} - \epsilon_{\vec{k}'} - \hbar\omega_0)] + (1/4)((k_y k'_z - k_z k'_y)^2 + (k_x k'_y - k_y k'_x)^2) \delta(\epsilon_{\vec{k}} - \epsilon_{\vec{k}'} + \hbar\omega_0)]$, where $\epsilon_{\vec{k}}$ is the energy of an electron with wave vector \vec{k} . The only effect occasioned by the passage to the rotating frame is the appearance of $\hbar\omega_0$ in the delta functions with quantum corrections to the form of the Gilbert term which are negligible since $\hbar\omega_0 \ll k_B T$. The modifications to the expression for $(1/\tau)$ are also negligible since, even for the true phononic mechanism, $\hbar\omega_0$ is much smaller than the relevant energy scale $k_B T$.

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- [1] J. C. Slonewski, J. Mag. Mag. Matr. **159**, L1 (1996).
 - [2] L. Berger. Phys. Rev. B **54**, 9353 (1996).
 - [3] J. J. Versluijs, M. A. Bari, and J. M. D. Coey Phys. Rev. Lett. **87**, 026601 (2001).
 - [4] D. A. Allwood, et al. Science **296**, 2003-2006 (2002).
 - [5] A. Yamaguchi et al., Phys. Rev. Lett. **92**, 077205 (2004).
 - [6] M. Yamanouchi, D. Chiba, F. Matsukura and H. Ohno, Nature **428**, 539 - 542 (2004)
 - [7] See, S. Zhang, P. M. Levy, and A. Fert, Phys. Rev. Lett. **88**, 236601 (2002); Y. Tserkovnyak, A. Brataas, G. E. W. Bauer, B. I. Halperin, cond-mat/0409242, and references therein.
 - [8] P. Matl et al., Phys. Rev. B **57**, 10248 (1998); Y. Taguchi et al., Science **291**, 2573 (2001); Y. Taguchi and Y. Tokura, Europhys. Lett. **54**, 401 (2001); S. H. Chun et al., Phys. Rev. Lett. **84**, 757 (2000).
 - [9] Y. B. Bazaliy, B. A. Jones and S.-C. Zhang Phys. Rev. B **57** R3213 (1998)
 - [10] G. Tataru, and H. Kohno, Phys. Rev. Lett. **92**, 086601 (2004).
 - [11] See e.g., C. Kittel, *Quantum theory of solids*, Wiley, New York (1963)
 - [12] When $K_\perp = 0$ there is a second rotational symmetry about the z -axis with S_{iz} , in the laboratory frame, as the local U(1) generator. It turns out that the single b_w^\dagger is also the Goldstone boson for this symmetry.
 - [13] One of the interests in the spintronics field is the possibility of exerting forces on, and pinning of, a domain wall. The presence of x' in the expression for J_s indicates that this potentially can be modified and modulated dynamically in, e.g., a field effect device.
 - [14] However $(1 - \cos\theta_i)(\nabla\phi_i)$ is a Berry phase which instead might be written as $\phi_i \sin\theta_i(\nabla\theta_i)$, implying an energy $\sim j_s\phi_i$ and a finite $\partial\mathcal{V}/\partial\phi_i$. These two fashions in which to write this phase *should* be related by a gauge transformation, i.e., $\vec{A} \rightarrow \vec{A} - \nabla f(\theta_i, \phi_i)$ where $f = (1 - \cos\theta_i)\phi_i\sigma_z$. This connection however fails since there is a large surface term at the end with $\theta_i = \pi$. The SU(2) rotations $r(\theta_i, \phi_i) \equiv e^{i\phi_i\sigma_z} e^{i\theta_i\sigma_y} e^{-i\phi_i\sigma_z}$ used here are *special* in that the gauge fields \vec{A} are zero far from the wall it creates and such surface terms are absent.
 - [15] See: Y. Yafet, Solid State Physics, Ed. F. Seitz and D. Turnbull (Academic, New York, 1963), **14**, 1.
 - [16] See review: S. E. Barnes, Adv. Phys. **30**, 801 (1981)